

Disulfur monoxide

Inchi: InChI=1S/OS2/c1-3-2
InchiKey: TXKMVPPZCYKFAC-UHFFFAOYSA-N
Formula: OS2
SMILES: O=S=S
Mol. weight [g/mol]: 80.13
CAS: 20901-21-7

Physical Properties

| Property code | Value | Unit | Source |
|---------------|--------------|----------------------|----------------|
| ea | 1.88 ± 0.01 | eV | NIST Webbook |
| gf | -130.71 | kJ/mol | Joback Method |
| hf | -95.58 | kJ/mol | Joback Method |
| hfus | 11.00 | kJ/mol | Joback Method |
| hvap | 34.84 | kJ/mol | Joback Method |
| ie | 10.62 | eV | NIST Webbook |
| ie | 10.58 ± 0.01 | eV | NIST Webbook |
| ie | 10.58 ± 0.01 | eV | NIST Webbook |
| ie | 10.30 ± 0.10 | eV | NIST Webbook |
| ie | 10.52 | eV | NIST Webbook |
| ie | 10.53 ± 0.02 | eV | NIST Webbook |
| ie | 10.58 ± 0.01 | eV | NIST Webbook |
| ie | 10.52 | eV | NIST Webbook |
| log10ws | 1.23 | | Crippen Method |
| logp | -0.339 | | Crippen Method |
| mvol | 45.130 | ml/mol | McGowan Method |
| pc | 9481.11 | kPa | Joback Method |
| tb | 325.06 | K | Joback Method |
| tc | 524.87 | K | Joback Method |
| tf | 193.38 | K | Joback Method |
| vc | 0.162 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-------|------|-----------------|--------|
|---------------|-------|------|-----------------|--------|

| | | | | |
|-----|-------|---------|--------|---------------|
| cpg | 52.17 | J/mol×K | 325.06 | Joback Method |
| cpg | 54.21 | J/mol×K | 358.36 | Joback Method |
| cpg | 55.91 | J/mol×K | 391.66 | Joback Method |
| cpg | 57.30 | J/mol×K | 424.97 | Joback Method |
| cpg | 58.40 | J/mol×K | 458.27 | Joback Method |
| cpg | 59.25 | J/mol×K | 491.57 | Joback Method |
| cpg | 59.87 | J/mol×K | 524.87 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C20901217&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| ea: | Electron affinity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| ie: | Ionization energy |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/48-616-5/Disulfur-monoxide.pdf>

Generated by Cheméo on 2024-04-17 01:49:40.860980605 +0000 UTC m=+15607829.781557921.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.